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Modified Formulation, ε -Regularization and the Efficient Solution of Hydraulic Fracture Problems

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Additional information is available at the end of the chapter

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1. Introduction

Analytical models and numerical simulation are important means to increase understanding and to enhance efficiency of hydraulic fracturing. The reasons for developing and using them are clearly explained, for instance, by Mack and Warpinski [1]. Thus, there is no need to dwell on them. Rather we focus on improving analytical and numeric methods used to the date. Our objective is to suggest new approaches for developing accurate, robust and stable simulators on the basis of recent analytical and computational findings [2-6].

The approaches discussed in the paper stem from the fact [2,3] that the conventional formulation of the hydraulic fracture problem (for example, see [7]), when neglecting the lag and fixing the position of the fracture front at a time step, is ill-posed. This feature has not been reported for more than three decades of studying hydraulic fractures because of two reasons. Numerical simulators, based on the conventional formulation (for example, see [7,8]), employ quite rough meshes, which themselves serve as specific 'regularizators'. On the other hand, rare solutions of model problems either also employed rough meshes [9], or they were obtained by solving the initial value (Cauchy) problem [10-12] rather than the boundary value problem (a discussion of the difference may be found in references [3,4]). The disclosure of the mentioned fact has led to (i) explicit formulation of the speed equation (SE) in its general form¹, (ii) comprehension of its significance for proper numerical simulation of hydraulic fractures and (iii) distinguishing the particle velocity as a preferable variable². It

¹ To the authors' knowledge, Kemp [11] was the first who clearly distinguished the speed equation when revisiting the Nordgren's problem. When introducing the SE, numbered (5) in his paper, Kemp wrote (p. 289): „Nowhere is (5) mentioned. (5) is called the Stefan condition and is always present in moving boundary problems“. Kemp used the SE to solve the Nordgren's problem as an initial value rather than boundary value problem. This excluded solving ill-posed boundary value problem.

has also led to the efficient means, called ε -regularization [2,3], to overcome the analytical and computational difficulties caused by ill-posedness. Finally, the entire conventional formulation of the problem has changed to the modified formulation, which opens new analytical and computational options for solving hydraulic fracture problems.

Below we employ these options. Section 2 contains a concise review of the modified formulation. In Section 3, we illustrate its analytical advantages by simple solutions for the Perkins-Kern-Nordgren (PKN) [9,13] and Khristianovich-Geertsma-de Klerk (KGD) [14,15] models. In Section 4, we turn to computational advantages of the modified formulation, present new computational schemes and illustrate their efficiency by numerical results for the PKN model. Section 5 contains the extension of these schemes to the pseudo-three-dimensional (P3D) models [1], the importance of which grows nowadays because of their employment in simulators for hydraulic fractures in low permeable shales [16]. A brief summary concludes the paper (Section 6).

2. Modified formulation of hydraulic fracture problem

The modified equations [3,4] use as variables the velocity v of fluid particles averaged through the channel width (fracture opening) and the modified opening $y=w^{1/\alpha}$, where α is the exponent defined by the asymptotic behavior of the opening w at the fluid front ($w=C(t)r^\alpha$, r is the distance from the front, t is the time). Then the continuity equation for a flow in a narrow channel reads:

$$\frac{\partial y}{\partial t} + \mathbf{v} \cdot \text{grad} y + \frac{1}{\alpha} y \text{div} \mathbf{v} + \frac{y^{1-\alpha}}{\alpha} q_l = 0 \quad (1)$$

where q_l is the intensity of distributed sinks or sources (below this term is assumed positive to account for leak off); the divergence and gradient are defined in the tangent plane to the channel (hydraulic fracture). The Poiseuille type relation in terms of the particle velocity is:

$$\mathbf{v} = -\frac{1}{w} D(w, p) \text{grad} p \quad (2)$$

It is obtained by integration of the Navier-Stokes equations for a flow of a viscous fluid in a narrow channel. Herein, $p(x, t)$ is the fluid pressure, x is the position vector on the surface of the flow in immovable coordinates, D is a function or operator, such that $D(0, p) \text{grad} p = 0$.

The equations (1) and (2) represent the modified system of PDE. Substitution of (2) into (1) gives the modified lubrication equation. We shall not write it down explicitly because in the

2 Mack and Warpinski [1] have noticed the beneficial property of the velocity. On p. 6-21 of their fundamental work they wrote: „Detailed numerical simulations have shown that the velocity varies much more slowly than the flow rate“. These authors made the best of this property by using the velocity as the unknown variable; they actually employed the speed equations, as well, although not writing it explicitly.

general case, keeping the velocity, which is a smooth function near the fluid front, as unknown is more convenient than to replace it by $\text{grad} p$, which is commonly strongly singular. Note only that the operator resulting from substitution of (2) into (1) is of first order in time and of second order and elliptic in spatial coordinates. Consequently it requires one *initial* condition defining the initial distribution of the opening. In terms of the modified opening it is:

$$y(x, t_0) = y_0(x) \quad (3)$$

The elliptic operator requires only one *boundary* condition (BC) on the fluid front L_e . For instance, when neglecting the lag between the fluid front and the fracture contour, the BC may be the condition of the prescribed normal component $q_n(x, t) = y^\alpha v_n$ of the flux:

$$y^\alpha v_n(x, t) = q_0(x, t), \quad x \in L_e \quad (4)$$

where $q_0(x)$ is a known function at L_e ; specifically, at the injection points, it is defined by the injection regime. At a point x_* of the propagating fluid front, coinciding with the fracture contour L_f , we have $w = 0$ and equation (2) implies $q_f(x, t) = wv = 0$. Then (4) becomes:

$$y(x_*, t) = 0, \quad x_* \in L_f \quad (5)$$

Of special importance is that the limit value of the particle velocity v_{n*} at the fluid front represents the speed of the front propagation v_* [2-4]:

$$v_* = \frac{dx_{n*}}{dt} = v_{n*} \quad (6)$$

Herein, x_{n*} is the normal component of a point x_* at the front. It is assumed that sucking or evaporation through the front is negligible. (6) is the *speed equation* (SE), which is fundamental for moving boundaries. Using (2) in (6) specifies the SE for a flow of incompressible fluid in a narrow channel [2,3]:

$$v_* = \frac{dx_{n*}}{dt} = - \left[\frac{1}{w(x_*)} D(w, p) \frac{\partial p}{\partial n} \right]_{x=x_*} \quad (7)$$

Thus we have the *local* condition (7) at points of the propagating fluid front. This allows one to trace the propagation by well-established methods of the theory of propagating surfaces [17]. In contrast, the conventional formulation employs the *global* mass balance (for example, see [7,8,10,12]), which is a single equation. The latter is sufficient when considering 1D problems with one point of the front to be traced. However, in the general case of a 2D fracture, it is preferable to employ the SE, which is formulated at each of the points of the fluid front.

This gives another evidence that using the particle velocity is beneficial from the computational point of view.

The speed equation (6) also yields important implications for numerical simulation of hydraulic fractures by finite differences. Indeed, when at a time step we have known both x_* and $v_*(x_*)$, the SE (7) becomes a boundary condition additional to the boundary condition (5) on the front. Thus, as noted in [2], a boundary value problem may appear overdetermined and ill-posed in the Hadamard sense [18]. To avoid difficulties, it is reasonable to use ε -regularization, suggested and employed in [2,3].

The ε -regularization is performed as follows. At each point of the fluid front, an exact BC is changed to an approximate equality at a small distance r_ε behind the front:

$$\int_{p_0}^{p_\varepsilon} \frac{1}{w} D(w, p) dp = v_* r_\varepsilon \quad (8)$$

where p_0 is the pressure at the front, p_ε is the pressure at a distance r_ε from the front. Equation (8) is obtained by combining the boundary condition at the fluid front, particular for a problem, with the SE, which is general. The distance (absolute r_ε or relative ε) is taken small enough to use the equality sign in the derived approximate condition. The SE is also assumed to be met at the distance r_ε to an accepted accuracy:

$$v_*(t) = \frac{dx_n}{dt} = - \left[\frac{1}{w} D(w, p) \frac{\partial p}{\partial n} \right]_{r=r_\varepsilon} \quad (9)$$

The ε -regularized BC (8) allows one to avoid unfavorable computational effects. The ε -regularized SE (9) serves to find the front propagation. In particular, it is basic when applying the level set and fast marching methods [17].

The described modification concerns mostly with the *fluid* equations, which are sufficient when the opening w is known. However, it is not known in advance in problems of hydraulic fracturing. To find it, the fluid equations are complimented with solid mechanics and fracture equations.

In terms of the modified opening, the *solid mechanics* equation with an integral operator A

$$Ay^\alpha = p \quad (10)$$

is solved under the BC of zero opening at points of the fracture contour. When neglecting the lag, this condition coincides with the condition (5) of zero flux at the fluid front.

The *fracture mechanics* equations define the critical state and the perspective direction of the fracture propagation. In the case of tensile mode of fracture, these are:

$$K_I = K_{Ic}, \quad K_{II} = 0 \quad (11)$$

where K_I is the tensile stress intensity factor (SIF), K_{Ic} is its critical value, K_{II} is the shear SIF. The modified problem, when neglecting the lag ($x_e = x_*$), consists of solving the PDE (1) and (2) together with the solid mechanics equation (10) under the initial condition (3), BC (4) at the contour part with prescribed influx, and the ϵ -regularized BC (8) at the fluid (and fracture) front. The conditions (11) define the possibility and the direction of the fracture propagation. The ϵ -regularized SE (9) serves to find the front propagation on the time steps.

3. Analytical solutions

In particular cases of 1D problems for the PKN [9,13] and KGD [14,15] models, one can solve an initial value rather than boundary value problem (for example, see [10-12]). Then the problem is well-posed and it does not require regularization. Still in these cases, the modified variables are of use to obtain simple analytical solutions of problems, which otherwise require involved calculations. For a Newtonian fluid, the analytical solutions are given in [4] both for the PKN and KGD models. In a similar way, by employing the modified variables, we may solve these 1D problems for a non-Newtonian fluid.

Consider a fluid with the viscosity law of power-type $\tau = M \dot{\gamma}^n$, where τ is the shear stress, $\dot{\gamma}$ is the shear strain rate, M is the consistency index, n is the behaviour index. The common derivations for a flow in a narrow channel yield the Poiseuille type dependence between the particle velocity v , averaged through the channel width w , and the pressure gradient:

$$v = \left(-k_f w^{n+1} \frac{\partial p}{\partial x} \right)^{1/n} \quad (12)$$

where $k_f = 1/(\theta M)$, for a plane channel, $\theta = 2[2(n+1)]^n$ (for example, see [12]), the x -axis is in the direction of the fracture propagation, v is the component of the particle velocity in the direction of x .

The geometrical scheme of the PKN model is given in Figure 1. In this model the plane-strain conditions occur in cross-sections parallel to the fracture front. Then the elasticity equation (10) takes the form [9]:

$$p = k_r w \quad (13)$$

where $k_r = [2/(\pi h)]E'$, $E' = E/(1-\nu^2)$, E is the elasticity modulus, ν is the Poisson's ratio, h is the fracture height. Substitution of (13) into (12) yields:

$$v = \left(-\frac{k_f k_r}{n+2} \frac{\partial w^{n+2}}{\partial x} \right)^{1/n} \quad (14)$$

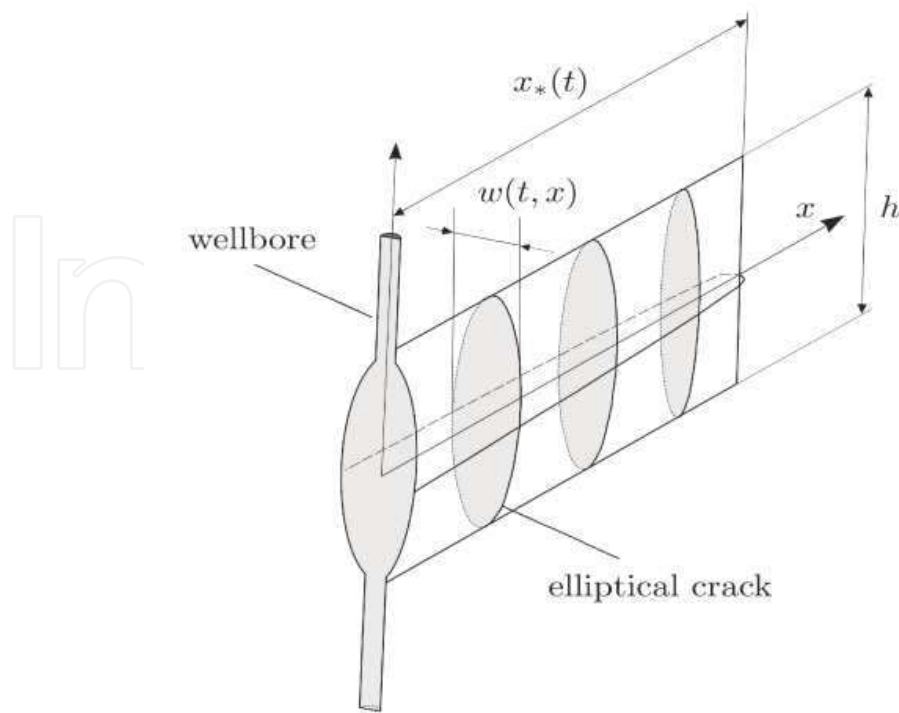


Figure 1. Geometrical scheme of the PKN model

Physically significant speed of the front is neither zero, nor infinite. According to (14) and the SE (7) it is possible only when the function $y = w^{n+2}$ is linear in x . Hence, in the considered problem, the exponent α in the modified opening $y = w^{1/\alpha}$ is $\alpha = 1 / (n + 2)$. Then the continuity equation (1), the velocity equation (2), the initial condition (3), the BC (4), (5) and the SE (6) become, respectively,

$$\frac{\partial y}{\partial t} + v \frac{\partial y}{\partial x} + \frac{y}{\alpha} \frac{\partial v}{\partial x} + \frac{y^{1-\alpha}}{\alpha} q_l = 0 \quad (15)$$

$$v = \left(-k_f k_r \alpha \frac{\partial y}{\partial x} \right)^{1/n} \quad (16)$$

$$y(x, t_0) = y_0(x) \quad (17)$$

$$[y^\alpha v]_{x=0} = q_0(t), \quad (18)$$

$$y(x_*) = 0 \quad (19)$$

$$v_*(t) = \frac{dx_*}{dt} = \left[\left(-k_f k_r \alpha \frac{\partial y}{\partial x} \right)^{1/n} \right]_{x=x_*} \quad (20)$$

Following [4], introduce the normalized variables $x_d = x / x_n$, $x_{*d} = x_* / x_n$, $t_d = t / t_n$, $v_d = v / v$, $v_{*d} = v_* / v_n$, $y_d = y / y_n$, $q_d = q / q_n$, $q_{ld} = q_l / q_{ln}$, where the normalizing quantities are: $x_n = (k_f k_r q_n^{n+2} t_n^{2n+2})^{1/(2n+3)}$, $v_n = x_n / t_n$, $y_n = (q_n t_n / x_n)^{1/\alpha}$ with t_n and q_n being arbitrary typical values of the time and influx per unit height, respectively. In terms of the normalized variables, the problem (15)-(20) has the same form except that the multiplier $k_f k_r$ is changed to the unit. This excludes the consistency factor, the elasticity modulus and the height. Hereafter in this section we use (15)-(20) assuming $k_f = 1$ and $k_r = 1$.

Consider the case when the influx is prescribed by the power dependence in time:

$$q_0(t) = t^{\beta_q} \quad (21)$$

, (21)

where β_q is a prescribed constant; for a constant influx, $\beta_q = 0$. In the case of zero leak-off, the solution of (15)-(20) with the influx (21) may be found in self-similar variables:

$$x = \xi t^{\beta_*} x_* = \xi_* t^{\beta_*} v = V(\xi) t^{\beta_*-1} v_* = V_* t^{\beta_*-1} y = Y(\xi) t^{\beta_w/\alpha} q = Y(\xi)^\alpha V(\xi) t^{\beta_q} \quad ()$$

with $\beta_w = [1 + (n+1)\beta_q] / (2n+3)$, $\beta_* = [2(n+1) + (n+2)\beta_q] / (2n+3)$. In (22), ξ_* and V_* are constants, expressing respectively the self-similar fracture length and the front speed.

We may account for leak-off by assuming that the term q_l is also represented in the form with separated variables: $q_l = Q_l(\xi) t^{\beta_l}$, where $Q_l(\xi)$ is a prescribed function, which may be singular at the front as $Q_l(\xi) = o((\xi_* - \xi)^{\alpha-1})$. For the exponent β_l , it follows that $\beta_l = \beta_w - 1$.

We prescribe the functions $Y(\xi)$ and $V(\xi)$ by power series in the variable $\tau = 1 - \xi / \xi_*$:

$$Y(\xi) = \frac{\xi_*^{n+1} \beta_*^n}{\alpha} \sum_{j=1}^{\infty} a_j \tau^j, \quad V(\xi) = V_* \sum_{j=0}^{\infty} b_j \tau^j, \quad (22)$$

where $V_* = \xi_* \beta_*$. The leak-off function $Q_l(\xi)$ is represented as $Q_l(\xi) = \tau^\alpha \sum_{j=0}^{\infty} q_j \tau^j$ with known coefficients q_j (for zero leak-off, $q_j = 0$, $j = 0, 1, \dots$). Then the coefficients b_j and a_j are found recurrently from the equations (15), (16) re-written in self-similar variables. Omitting technical details, the coefficients b_j for $j = 2, 3, \dots$ are:

$$b_j = -\frac{1}{j+\alpha} \left\{ \sum_{k=2}^j \left[(j-k+1+\alpha k) a_k b_{j-k+1} + \left(\alpha j - \frac{\beta_w}{\beta_*} \right) a_j \right] - C_l \sum_{k=1}^j c_k q_{j-k} \right\} \quad (23)$$

with the starting values $a_1=b_0=1$, $b_1=\frac{1}{1+\alpha}\left(-\alpha+\frac{\beta_w}{\beta_*}+C_l q_0\right)$, $c_1=1$. In (24), $C_l=\left(\frac{\alpha}{\xi_*^{n+1}\rho_*^{n+1/\alpha}}\right)^\alpha$ and the coefficients a_j and c_j are found recurrently from equations: $\sum_{k=0}^{\infty}(k+1)a_{k+1}\tau^k=\left(\sum_{j=0}^{\infty}b_j\tau^j\right)^n$, $\sum_{k=1}^{\infty}c_k\tau^k=\tau\left(\sum_{j=0}^{\infty}a_{j+1}\tau^j\right)^\alpha$. In particular, for the first five coefficients a_j and c_j we have:

$$\begin{aligned} a_1=b_0=1, \quad a_2=\frac{1}{2}nb_1, \quad a_3=\frac{1}{6}n[(n-1)b_1^2+2b_2] \\ a_4=\frac{1}{24}n[(n-1)(n-2)b_1^3+6(n-1)b_1b_2+6b_3] \end{aligned} \quad (24)$$

$$a_5=\frac{1}{120}n[(n-1)(n-2)(n-3)b_1^4+12(n-1)(n-2)b_1^2b_2+24(n-1)b_1b_3+24b_4] \quad (25)$$

$$\begin{aligned} c_1=a_1=1, \quad c_2=(1-\alpha)a_2, \quad c_3=\frac{1}{2}(1-\alpha)[- \alpha a_2^2+2a_3], \quad c_4=\frac{1}{6}(1-\alpha)[\alpha(\alpha+1)a_2^3-6\alpha a_2a_3+6a_4], \\ c_5=\frac{1}{24}(1-\alpha)[- \alpha(\alpha+1)(\alpha+2)a_2^4+12\alpha(\alpha+1)a_2^2a_3-24\alpha a_2a_4+24a_5]. \end{aligned} \quad (26)$$

Starting from $a_1=b_0=c_1=1$, $b_1=(-\alpha+\beta_w/\beta_*+C_l q_0)/(1+\alpha)$, we find a_2 from the second of (25) and c_2 from the second of (26). Then (24) gives b_2 , the third of (25) defines a_3 , the third of (26) defines c_3 and so on.

The value ξ_* of the self-similar fracture length is found from the self-similar BC at the inlet: $[Y(0)]^\alpha V(0)=1$. By using the solution (23)-(26) for various ξ_* , we find the one, which meets the BC with a prescribed tolerance.

In the case of a *Newtonian fluid* ($n=1$, $\alpha=1/3$) equations (23)-(26) extend the analytical solution, obtained in [4], to the case of non-zero leak-off. In the case of a *perfectly plastic fluid* ($n=0$, $\alpha=1/2$), we have $a_k=0$, $c_k=0$ for $k>1$. Then for constant influx ($\beta_q=0$), the solution is:

$\xi_*= (9/8)^{1/3}$, $Y(\xi)=2(\xi_*-\xi)$, $V(\xi)=V_*\left[1+\frac{1}{\beta_*\sqrt{2\xi_*}}\sum_{j=1}^{\infty}\frac{2}{2j+1}q_{j-1}\tau^j\right]$, $V_*=2/3\xi_*$. Calculations show that the difference between the self-similar solutions for these two limiting cases is quite small. For instance, for zero leak-off, the self-similar fracture length is $\xi_{*p}=1.04004$ for a perfectly plastic fluid; it is $\xi_{*N}=1.00101$ for a Newtonian fluid. In both cases, the particle velocity is almost constant along the fracture. Thinning fluids ($0<n<1$), being intermediate between Newtonian and perfectly plastic fluids, the conclusions hold for an arbitrary thinning fluid. A detailed discussion of the solution is given in [19].

4. Increasing efficiency of numerical simulations

As mentioned, in the general case, using the SE (7) opens options for tracing the fracture in the 3D space by level set, fast marching and other methods of the theory of propagating sur-

faces [17]. As to the authors' knowledge, to this date, the only paper, in which the level set method has been applied to the hydraulic fractures, is that by Peirce and Detournay [20]. Since these authors used the conventional formulation not including the SE, special technique, called by them "implicit level set method", was suggested and successfully used. Still, direct employing of the SE looks superior in simplicity and possibility to use the standard well-established technique.

From now on, we focus on other computational advantages of the modified formulation which evidently appear when considering the PKN model. Recall that this model is basic for a wide class of simulators employing the P3D models (for example, see [1]). Since the Nordgren's problem is 1D, it is convenient to use the spatial coordinate x normalized by the fracture length x_* : $\varsigma = x / x_*$. In this spatial variable, the problem (15)-(20) reads:

$$\frac{\partial y}{\partial t} = -\frac{y}{\alpha x_*} \frac{\partial v}{\partial \varsigma} + \frac{\varsigma v_* - v}{x_*} \frac{\partial y}{\partial \varsigma} - \frac{y^{1-\alpha}}{\alpha} q_l, \quad (27)$$

$$v = \frac{k}{x_*^{1/n}} \left(-\frac{\partial y}{\partial \varsigma} \right)^{1/n} \quad (28)$$

$$y(\varsigma, t_0) = y_0(\varsigma) \quad (29)$$

$$[y^\alpha v]_{\varsigma=0} = q_0(t), \quad (30)$$

$$y(1, t) = 0 \quad (31)$$

$$v_*(t) = \frac{d x_*}{dt} = \frac{k}{x_*^{1/n}} \left[\left(-\frac{\partial y}{\partial \varsigma} \right)^{1/n} \right]_{\varsigma=1} \quad (32)$$

where $k = (k_f k_r \alpha)^{1/n}$. When writing (27)-(32), we have used the symbols y and v for the functions $\tilde{y}(\varsigma, t) = y(x(\varsigma), t)$ and $\tilde{v}(\varsigma, t) = v(x(\varsigma), t)$ omitting the tilde. Note that the initial condition (28) defines also the initial value x_{*0} of the fracture length as the end point where the opening is zero. Thus we actually have the initial condition for the length:

$$x_*(t_0) = x_{*0} \quad (33)$$

In the case when $y_0(\varsigma) = 0$, we have $x_{*0} = 0$.

In the considered problem the dependence (28) between the unknowns is explicit. Therefore, we may substitute (28) into (27). This yields the PDE of the first order in time and of the second order in ς :

$$\frac{\partial y}{\partial t} = \frac{k}{z_*} \left\{ \frac{y}{\alpha n} \left(-\frac{\partial y}{\partial \zeta} \right)^{(1-n)/n} \frac{\partial^2 y}{\partial \zeta^2} + \left[\zeta \left(-\frac{\partial y}{\partial \zeta} \right)^{1/n} - \left(-\frac{\partial y}{\partial \zeta} \right)^{1/n} \right] \frac{\partial y}{\partial \zeta} \right\} - \frac{y^{1-\alpha}}{\alpha} q_l \quad (34)$$

where $z_* = x_*^{(1+n)/n}$ is the modified fracture length. For it, the SE (32) becomes:

$$\frac{dz_*}{dt} = \frac{n+1}{n} k \left[\left(-\frac{\partial y}{\partial \zeta} \right)^{1/n} \right]_{\zeta=1} \quad (35)$$

From (34) and (35) it follows that after spatial discretization of these equations and the BC (30), (31), we obtain a well-posed system of ordinary differential equations (ODE) of the first order in time to be solved under the initial conditions (29) and (33). Actually, the problem does not require regularization. Still, as noted in [6], not to have too stiff system of ODE, it is reasonable to employ the ε -regularized forms of the BC (31) and the SE (34):

$$y(1-\varepsilon, t) = \left[\frac{n}{k(n+1)} \frac{dz_*}{dt} \right]^n \varepsilon \quad (36)$$

$$\frac{dz_*}{dt} = \frac{n+1}{n} k \left[\left(-\frac{\partial y}{\partial \zeta} \right)^{1/n} \right]_{\zeta=1-\varepsilon} \quad (37)$$

where ε is a small positive value in the range from 10^{-3} to 10^{-6} (it may yet be less depending on round-off errors of computer calculations).

Solving the system of ODE, resulting from spatial discretization of (34), (35), under the initial conditions (30), (33) may be performed by well-developed methods, like the Runge-Kutta method. Standard solvers are of immediate use. Emphasize that this option has appeared only due to employing the *local* SE rather than the global mass balance for tracing the front propagation. As show numerous numerical experiments, summarized in [6] for a Newtonian fluid ($n = 1$), this computational scheme provides highly accurate and stable results with small time expense.

As an example, we use this scheme to extend the Nordgren's numerical results [9] on the dependence of the fracture length on time for a Newtonian fluid and Carter's leak-off. To this end, the normalizing length x_n and normalizing time t_n are taken as those in the paper by Nordgren:

$$x_n = \pi \left(\frac{\mu q_0^5 h}{128 E C^8} \right)^{1/3}, \quad t_n = \pi^2 \left(\frac{\mu q_0^2 h}{16 E C^5} \right)^{2/3}.$$

Herein, μ is the dynamic viscosity, C is the fluid-loss coefficient in the Carter's leak-off term $q_l = 2C / \sqrt{t - \tau}$, τ is the time at which the fracture front reaches a point x . In accordance with (18), we have used the influx q_0 per unit height, while Nordgren wrote his normalizing values in terms of the total influx q_1 through the entire height h ($q_1 = q_0 h$).

The results are summarized in Figure 2. The thick solid line presents the Nordgren's curve (Fig. 2 of reference [9]), obtained in the time range [0.01, 5.0]. The author did not comment on the accuracy of his calculations. The accuracy, as stated in [3], is certainly below 1% which is also evident from the fact that the Nordgren's curve intersects the asymptotic dotted line, corresponding to small time, and the asymptotic dashed line, corresponding to large time.

The results obtained by employing the computational scheme described are presented in Figure 2 by the thin solid line. It can be seen that the solution starting from the small time asymptotics tends to that corresponding to the large time one. The calculations are performed by using ϵ -regularized BC (36) and SE (37) with $\epsilon=10^{-4}$. Specially designed numerical experiments have shown that the relative error does not exceed 10^{-3} . No signs of instability are noted. Even though the calculations were performed in the MATLAB environment with use of the standard ODE solver *ode15s*, the run time to cover the time range of 12 orders is 10 seconds. The said confirms high efficiency of the computational scheme suggested.³

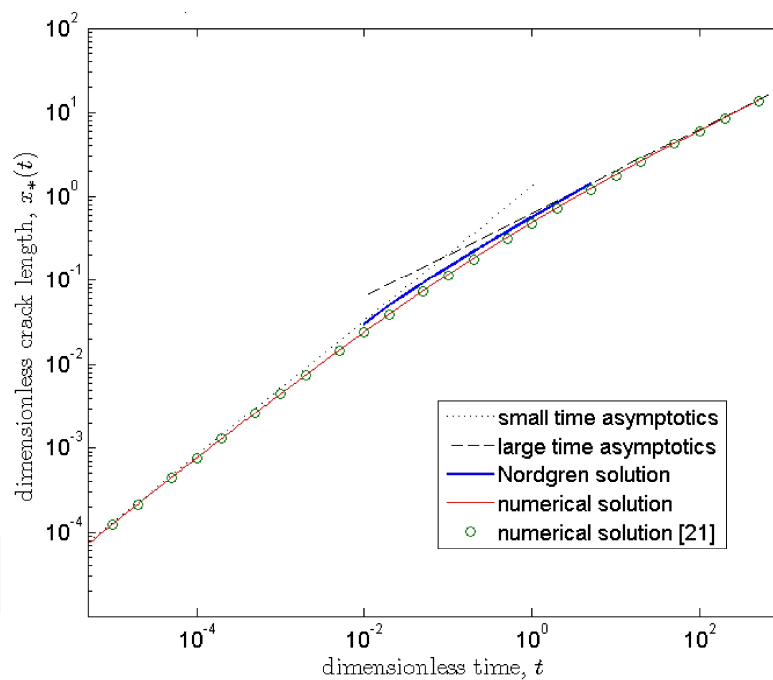


Figure 2. Dimensionless fracture length vs dimensionless time

³ Previously, Kovalyshen and Detournay [21] also have solved the same problem in the range of time wider than that covered by Nordgren. The authors performed calculations starting from the small-time asymptote as an initial condition at $\tau=10^{-8}$ and presented numerical results in the range $10^{-5} < \tau < 500$. Their results are shown by circles in Figure 2. They are indistinguishable from our numerical solution. The authors do not discuss the accuracy, stability and robustness of their calculations. From our results, to which the relative error does not exceed 10^{-3} , we may conclude that the relative error of the results, given for the fracture length in Table 1 of the paper [21], does not exceed 1% in the whole range of the calculations.

A similarly efficient computational scheme consists of solving the PDE (27), (28) with ε -regularized BC and SE by the Crank-Nicolson method [3,6].

Comment. In view of the accepted linear dependence (13) between the net-pressure and the opening, the pressure may replace the opening as an unknown variable. Actually, under the normalization, yielding $k_r=1$, there is no difference between p and w .

5. Extension to P3D models

In the previous section we have stated that the modified formulation provides prerequisites for efficient solving the Nordgren's problem. The latter, being the basis for the P3D models, we may extend the efficient schemes to these models.

The detailed description of the P3D models is given by Mack and Warpinski [1]. Thus it is sufficient to list only those their features, which distinguish them from the PKN model (Fig. 1) and which are significant for the extension.

(i) In P3D models, the in-situ rock stresses are different in various layers while the fluid pressure is assumed constant in any vertical cross-section. This implies that, in contrast with the PKN model, the net-pressure is now not constant in a vertical cross-section. Thus the P3D models employ the pressure itself rather than the net-pressure. Alternatively, one may employ the difference of the fluid pressure with a *fixed* value of in-situ rock pressure, say the pressure in the pay-layer. Below to keep clear connection with the PKN model, we shall use this option and conditionally call the difference the 'net-pressure'.

(ii) The linear dependence (13) between the net-pressure p and the average opening w_{av} of a cross-section is changed to a non-linear dependence $p=p(w_{av})$. To have clear connection with the initial PKN model, we re-write it as

$$p=k_r w_{av} F_p(w_{av}) \quad (38)$$

where for sufficiently small w_{av} , in particular near the fracture front, $F_p(w_{av})=1$. A specific form of the dependence (38) is found from solving plane-strain elasticity equation for a straight vertical crack under the fracture conditions of the form $K_I \leq K_{Ic}$ at the upper and bottom tips. Although looking for this dependence may be involved, it is found in advance for a prescribed geometry of layered stratum, in-situ stresses in layers and critical SIFs. These preliminary calculations also give the positions $z_u(p)$ and $z_l(p)$ of the upper and bottom tips, respectively. Consequently, the height $h_f(p)$ of the fracture in a cross section is a known function of the pressure:

$$h_f(p)=z_u(p)-z_l(p) \quad (39)$$

(iii) The dependence (12) between the particle velocity and the pressure, after averaging the velocity over the height of a cross-section, obtains a factor $F_v(w_{av})$:

$$v_{av} = \left(-k_f w_{av}^{n+1} \frac{\partial p}{\partial x} \right)^{1/n} F_v(w_{av}) \quad (40)$$

where for sufficiently small w_{av} , we have $F_v(w_{av})=1$. Since $w_{av} \rightarrow 0$ when approaching the fracture front, the SE becomes:

$$v_* = \left(-k_f w_{av}^{n+1} \frac{\partial p}{\partial x} \right)^{1/n}_{x=x_*} \quad (41)$$

(iv) The continuity equation is integrated over the cross-sectional height. As a result, the total flux through a cross-section and the total leak-off replace, respectively, the flux and leak-off per unit height. The average velocity v_{av} , being defined as the ratio of the total flux to the area A of the cross-section, the area A replaces the opening present in the PKN model. To preserve connection with the PKN model, we may use the flux, leak-off losses and area divided by a fixed reference height h_r , say the height of the pay-layer. Then denoting

$$w = \frac{A}{h_r}, \quad q = \frac{Av_{av}}{h_r} = wv_{av}, \quad q_l = \frac{Q_l}{h_r}, \quad (42)$$

we have the modified continuity equation (15) in the unchanged form. The BC (18) at the inlet is also the same when denoting $q_0(t) = Q_0(t)/h_r$, where $Q_0(t)$ is the prescribed total influx.

Take into account that by the definition of the average opening, we have $A = w_{av} h_f$, where h_f is the fracture height in a considered cross-section. Thus, by the first of (42), $w = Ah_f / h_r$. Then, in view of (39) and (38), we may use the argument w instead of w_{av} in the equations (38), (40) and (41) writing them, respectively, as

$$p = k_r w G_p(w), \quad v = \left(-k_f w^{n+1} \frac{\partial p}{\partial x} \right)^{1/n} G_v(w), \quad \text{and} \quad v_* = \left(-k_f w^{n+1} \frac{\partial p}{\partial x} \right)^{1/n}_{x=x_*}$$

where to simplify notation, we have omitted the subscript in the averaged velocity v_{av} . The functions $G_p(w)$ and $G_v(w)$ are found in advance through the functions $F_p(w_{av})$ and $F_v(w_{av})$. They are such that $G_p(w)=1$ and $G_v(w)=1$ for sufficiently small w .

Finally, by introducing the variables $y = w^{1/\alpha}$ and $\zeta = x / x_*$, we arrive at the system similar to (27)-(33):

$$\frac{\partial y}{\partial t} = -\frac{y}{\alpha x_*} \frac{\partial v}{\partial \zeta} + \frac{\zeta v_* - v}{x_*} \frac{\partial y}{\partial \zeta} - \frac{y^{1-\alpha}}{\alpha} q_l, \quad (43)$$

$$v = \frac{k}{x_*^{1/n}} H(y) \left(-\frac{\partial y}{\partial \varsigma} \right)^{1/n} \quad (44)$$

$$y(\varsigma, t_0) = y_0(\varsigma) \quad (45)$$

$$[y^\alpha v]_{\varsigma=0} = q_0(t), \quad (46)$$

$$y(1, t) = 0 \quad (47)$$

$$v_*(t) = \frac{dx_*}{dt} = \frac{k}{x_*^{1/n}} \left[\left(-\frac{\partial y}{\partial \varsigma} \right)^{1/n} \right]_{\varsigma=1} \quad (48)$$

$$x_*(t_0) = x_{*0} \quad (49)$$

where

$$H(y) = \left[\frac{d(wG_p(w))}{dw} \right]^{1/n} G_v(w) \quad (50)$$

is a function to be found in advance through the functions $G_p(w)$ and $G_v(w)$ of the argument $w = y^\alpha$. By the properties of $G_p(w)$ and $G_v(w)$, we have $H(0) = 1$ which explains its absence in the SE (48).

The problem (43)-(49) differs from the problem (27)-(33) in the only detail: equation (44) for the velocity contains the function $H(y)$ defined by (50). The latter function, being smooth and tending to the unity at the front, the efficient numerical schemes, discussed in the previous section, are of use for the P3D models.

Comment. In some cases, it may be convenient to use the net-pressure rather than the opening. Since $dp = \left(\frac{dp}{dw} \right) dw$, reformulation of the equations and computational schemes in terms of the modified pressure $P = p^{1/\alpha}$ is obvious.

6. Conclusions

The discussion above demonstrates the analytical and computational advantages of using the modified formulation. The analytical advantages are evident from the obtained simple analytical solutions for the PKN and KGD models, which otherwise require involved calculations. The computational advantages include: (i) the possibility to use the well-established theory of propagating surfaces, (ii) avoiding deterioration of numerical solution caused by ill-posedness of the problem when neglecting the lag and fixing the fracture contour at a

time step, (iii) avoiding singularities on the fluid front, (iv) the possibility to use highly efficient numerical schemes for the PKN and P3D models. These beneficial features are of significance when developing simulators able to efficiently solve truly 3D and pseudo-three-dimensional problems in real time. The work on simulators of both types is in progress.

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